

=> fil hcaplus
FILE 'HCAPLUS' ENTERED AT 10:01:27 ON 10 FEB 2003
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FILE COVERS 1907 - 10 Feb 2003 VOL 138 ISS 7
FILE LAST UPDATED: 7 Feb 2003 (20030207/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=>
=>

=> d stat que
L1 1 SEA FILE=REGISTRY ABB=ON PLU=ON PVLDLLRELLEELKQKLK/SQSP
L2 2 SEA FILE=HCAPLUS ABB=ON PLU=ON L1

=>
=>

=> d ibib abs hitrn l2 1-2

L2 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2003 ACS
ACCESSION NUMBER: 1999:233772 HCAPLUS
DOCUMENT NUMBER: 130:262129
TITLE: Apolipoprotein A-I .alpha.-helical peptide analogs as agonists for treatment of dyslipidemias
INVENTOR(S): Dasseux, Jean-Louis; Sekul, Renate; Buttner, Klaus; Cornut, Isabelle; Metz, Gunther; Dufourcq, Jean
PATENT ASSIGNEE(S): USA
SOURCE: PCT Int. Appl., 232 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9916409	A2	19990408	WO 1998-US20329	19980928
W:	AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, PG, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM			
RW:	CH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MF, NE, SN, TD, TG			

CA 2304814 AA 19990408 CA 1998-2304814 19980928
EP 1039934 A1 20001004 EP 1998-950742 19980928
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
IE, FI
NO 2000001601 A 20000526 NO 2000-1601 20000328
PRIORITY APPLN. INFO.: US 1997-940136 A 19970929
WO 1998-US20329 W 19980928

OTHER SOURCE(S): MARPAT 130:262129

AB Analogs of the .alpha.-helical peptides of apolipoprotein A-I (ApoA-I) that can act as ApoA-I agonists or superagonists with many at least as active as wild-type ApoA-I are described for use in treatment of dyslipidemias. Genes for these peptides may be used in gene therapy (no data). Detail physicochem. requirements for the amphipathic .alpha.-helices are given and these are quite different from the prior art understanding of the properties of amphipathic .alpha.-helices of ApoA-I. A series of >250 amphipathic peptides were tested for their ability to activate LCAT. One of these peptides was found to stimulate the formation of HDL with incorporation of cholesterol.

IT 221884-94-2

RL: BPR (Biological process); BSU (Biological study, unclassified); PRP (Properties); THU (Therapeutic use); BIOL (Biological study); PROC (Process); USES (Uses)
(activation of LCAT by, as amphipathic .alpha.-helical ApoAI agonist; apolipoprotein A-I .alpha.-helical peptide analogs as agonists for treatment of dyslipidemias)

L2 ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2003 ACS

ACCESSION NUMBER: 1999:233771 HCAPLUS

DOCUMENT NUMBER: 130:262128

TITLE: Apolipoprotein A-I agonists and their use to treat dyslipidemic disorders

INVENTOR(S): Dasseux, Jean-Louis; Sekul, Renate; Buttner, Klaus; Cornut, Isabelle; Metz, Gunther

PATENT ASSIGNEE(S): USA

SOURCE: PCT Int. Appl., 245 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9916408	A2	19990408	WO 1998-US20328	19980928
W:	AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, PQ, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TC, TM			
RW:	GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG			
US 6037323	A	20000314	US 1997-940093	19970929
CA 2304931	AA	19990408	CA 1998-2304931	19980928
EP 1019010	A1	20000719	EP 1998-951979	19980928
R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, FI			
AU 747823	B2	20020523	AU 1998-97791	19980928
US 6265377	B1	20010724	US 1999-465719	19991217
NO 2000001600	A	20000503	NO 2000-1600	20000328
US 2003008827	A1	20030109	US 2001-865989	20010325
PRIORITY APPLN. INFO.:			US 1997-940093 A 19970929	
			WO 1998-US20328 W 19980928	

OTHER SOURCE(S): MARPAT 130:262128

AB The present invention provides peptides and peptide analogs that mimic the structural and pharmacol. properties of human ApoA-I. The peptides and peptide analogs are useful to treat a variety of disorders assocd. with dyslipidemia.

IT 221884-94-2P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); PNU (Preparation, unclassified); PRP (Properties); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(apolipoprotein A-I agonists and their use to treat dyslipidemic disorders)

=>

=>

=> fil reg

FILE 'REGISTRY' ENTERED AT 10:01:38 ON 10 FEB 2003

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 7 FEB 2003 HIGHEST RN 487578-67-6

DICTIONARY FILE UPDATES: 7 FEB 2003 HIGHEST RN 487578-67-6

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=> d .seq 11 1

L1 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS

RN 221884-94-2 REGISTRY

CN L-Lysinamide, 1-acetyl-L-prolyl-L-valyl-L-leucyl-L-.alpha.-aspartyl-L-leucyl-L-leucyl-L-arginyl-L-.alpha.-glutamyl-L-leucyl-L-leucyl-L-.alpha.-glutamyl-L-.alpha.-glutamyl-L-leucyl-L-lysyl-L-glutamyl-L-lysyl-L-leucyl-
 (9CI) (CA INDEX NAME)

NTE modified

type	location		description
terminal mod.	Pro-1	-	N-acetyl
terminal mod.	Lys-18	-	C-terminal amide

SQL 18

SEQ 1 PVLDLLRELL EELKQKLLK

=====

HITS AT: 1-18

REFERENCE 1: 130:262129

REFERENCE 2: 130:262128

OM protein - protein search, using sw model

Run on: February 10, 2003, 09:49:05 ; Search time 29 Seconds
(without alignments)
127.891 Million cell updates/sec

Title: US-09-865-989-191
Perfect score: 85
Sequence: 1 PVLDLLRELLEELKQKLK 18

Scoring table: BLOSUM62
Gapop 10.0 , Gapext 0.5

Searched: 671580 seqs, 206047115 residues

Total number of hits satisfying chosen parameters: 671580

Minimum DB seq length: 0
Maximum DB seq length: 2000000000

Post-processing: Minimum Match 0%
Maximum Match 100%
Listing first 45 summaries

Database : SPTREMBL_21:*

- 1 sp_archaea:*
- 2 sp_bacteria:*
- 3 sp_fungi:*
- 4 sp_human:*
- 5 sp_invertebrate:*
- 6 sp_mammal:*
- 7 sp_mhc:*
- 8 sp_organelle:*
- 9 sp_phage:*
- 10: sp_plant:*
- 11: sp_rodent:*
- 12: sp_virus:*
- 13: sp_vertebrate:*
- 14: sp_unclassified:*
- 15: sp_rvirus:*

16: sp_bacteriap.*
17: sp_archeap.*

Pred. No. is the number of results predicted by chance to have a score greater than or equal to the score of the result being printed, and is derived by analysis of the total score distribution.

SUMMARIES

		%					
Result	Query						
No.	Score	Match	Length	DB	ID	Description	
1	49.5	58.2	1156	17	Q8TK58	Q8tk58 methanosarc	
2	49	57.6	182	16	Q9WXZ8	Q9wxz8 thermotoga	
3	47	55.3	281	10	Q9FGU2	Q9fgu2 arabidopsis	
4	47	55.3	658	5	Q9U3N1	Q9u3n1 caenorhabdi	
5	46	54.1	383	17	Q8U0V9	Q8u0v9 pyrococcus	
6	46	54.1	395	17	Q8U430	Q8u430 pyrococcus	
7	46	54.1	413	5	Q25004	Q25004 homarus ame	
8	45	52.9	277	16	Q92Q82	Q92q82 rhizobium m	
9	45	52.9	341	16	Q9RVB3	Q9rvb3 deinococcus	
10	45	52.9	447	16	Q8RD37	Q8rd37 thermoanaer	
11	45	52.9	488	5	Q9U3E8	Q9u3e8 caenorhabdi	
12	45	52.9	999	2	Q93JY2	Q93jy2 erwinia chr	
13	44.5	52.4	84	17	Q8TYE1	Q8tye1 methanopyru	
14	44	51.8	96	15	P89843	P89843 human immun	
15	44	51.8	135	2	Q93IB0	Q93ib0 staphylococ	
16	44	51.8	327	16	Q8R794	Q8r794 thermoanaer	
17	44	51.8	334	16	Q9A8K6	Q9a8k6 caulobacter	
18	44	51.8	343	2	P77840	P77840 chloroflexu	
19	44	51.8	364	5	Q17983	Q17983 caenorhabdi	
20	44	51.8	398	4	Q96KY5	Q96ky5 homo sapien	
21	44	51.8	458	16	Q8R8N9	Q8r8n9 thermoanaer	
22	44	51.8	487	16	Q98QF4	Q98qf4 mycoplasma	
23	44	51.8	566	16	Q9PHU2	Q9phu2 campylobact	
24	44	51.8	585	4	Q8WW52	Q8ww52 homo sapien	
25	44	51.8	791	17	Q9YB89	Q9yb89 aeropyrum p	
26	44	51.8	842	16	Q92FU6	Q92fu6 listeria in	
27	44	51.8	842	16	Q8YAV6	Q8yav6 listeria mo	
28	44	51.8	1078	5	Q18476	Q18476 caenorhabdi	
29	44	51.8	1942	17	Q8TVI4	Q8tvi4 methanopyru	
30	43.5	51.2	216	17	O58396	O58396 pyrococcus	
31	43	50.6	199	15	Q9JCT2	Q9jct2 human immun	
32	43	50.6	203	15	Q9JCX0	Q9jcx0 human immun	
33	43	50.6	341	16	Q8YUZ2	Q8yuz2 anabaena sp	

34	43	50.6	379	3	Q9P635	Q9p635 neurospora
35	43	50.6	383	17	Q9V0F0	Q9v0f0 pyrococcus
36	43	50.6	396	3	Q9USQ4	Q9usq4 schizosacch
37	43	50.6	398	17	O58030	O58030 pyrococcus
38	43	50.6	398	17	Q9UY32	Q9uy32 pyrococcus
39	43	50.6	405	2	Q04389	Q04389 bacillus sp
40	43	50.6	426	16	Q929T3	Q929t3 listeria in
41	43	50.6	462	12	Q9IBR4	Q9ibr4 spodoptera
42	43	50.6	469	16	Q8R8Q4	Q8r8q4 thermoanaer
43	43	50.6	522	12	Q9YVT0	Q9yvt0 melanoplus
44	43	50.6	571	17	Q980V7	Q980v7 sulfolobus
45	43	50.6	612	10	O65437	O65437 arabidopsis

ALIGNMENTS

RESULT 1

Q8TK58

ID Q8TK58 PRELIMINARY; PRT; 1156 AA.

AC Q8TK58;

DT 01-JUN-2002 (TrEMBLrel. 21, Created)

DT 01-JUN-2002 (TrEMBLrel. 21, Last sequence update)

DT 01-JUN-2002 (TrEMBLrel. 21, Last annotation update)

DE Phosphorylase.

GN MA3560.

OS Methanosarcina acetivorans.

OC Archaea; Euryarchaeota; Methanococci; Methanosarcinales;

OC Methanosarcinaceae; Methanosarcina.

OX NCBI_TaxID=2214;

RN [1]

RP SEQUENCE FROM N.A.

RC STRAIN=C2A / ATCC 35395 / DSM 2834;

RX MEDLINE=21929760; PubMed=11932238;

RA Galagan J.E., Nusbaum C., Roy A., Endrizzi M.G., Macdonald P.,

RA FitzHugh W., Calvo S., Engels R., Smirnov S., Atnoor D., Brown A.,

RA Allen N., Naylor J., Stange-Thomann N., DeArellano K., Johnson R.,

RA Linton L., McEwan P., McKernan K., Talamas J., Tirrell A., Ye W.,

RA Zimmer A., Barber R.D., Cann I., Graham D.E., Grahame D.A., Guss A.M.,

RA Hedderich R., Ingram-Smith C., Kuettner H.C., Krzycki J.A.,

RA Leigh J.A., Li W., Liu J., Mukhopadhyay B., Reeve J.N., Smith K.,

RA Springer T.A., Umayam L.A., White O., White R.H., de Macario E.C.,

RA Ferry J.G., Jarrell K.F., Jing H., Macario A.J.L., Paulsen I.,

RA Pritchett M., Sowers K.R., Swanson R.V., Zinder S.H., Lander E.,

RA Metcalf W.W., Birren B.;

RT "The genome of Methanosarcina acetivorans reveals extensive metabolic
RT and physiological diversity.";
RL Genome Res. 12:532-542(2002).
DR EMBL; AE011064; AAM06921.1; -.
KW Complete proteome.
SQ SEQUENCE 1156 AA; 130192 MW; 0280A362DC4BB5C6 CRC64;

Query Match 58.2%, Score 49.5; DB 17; Length 1156;
Best Local Similarity 57.9%, Pred. No. 1.7e+02;
Matches 11; Conservative 5; Mismatches 2; Indels 1; Gaps 1;

Qy 1 PVLDL-LRELLEELKQK 18
|:| |::|||::|
Db 749 PILDRNLKELLEIQKKAK 767

Search completed: February 10, 2003, 09:51:07
Job time : 32 secs

OM protein - protein search, using sw model

Run on: February 10, 2003, 09:49:05 ; Search time 11 Seconds
(without alignments)
67.870 Million cell updates/sec

Title: US-09-865-989-191
Perfect score: 85
Sequence: 1 PVLDLLRELLEELKQK 18

Scoring table: BLOSUM62
Gapop 10.0 , Gapext 0.5

Searched: 112892 seqs, 41476328 residues

Total number of hits satisfying chosen parameters: 112892

Minimum DB seq length: 0
Maximum DB seq length: 2000000000

Post-processing: Minimum Match 0%
Maximum Match 100%
Listing first 45 summaries

Database : SwissProt_40:*

Pred. No. is the number of results predicted by chance to have a score greater than or equal to the score of the result being printed, and is derived by analysis of the total score distribution.

SUMMARIES

Result No.	% Query					Description
	Score	Match	Length	DB	ID	
1	49	57.6	378	1	2NPD_NEUCR	Q01284 neurospora
2	47	55.3	439	1	MYC_MARMO	P22555 marmota mon
3	47	55.3	785	1	PTA1_YEAST	Q01329 saccharomyc
4	47	55.3	833	1	GYRA_BACHD	O50628 bacillus ha
5	46	54.1	1453	1	Y373_BOVIN	Q9tu23 bos taurus

6	45	52.9	558	1	YCXA_ASTLO
7	44	51.8	134	1	YTKC_BACSU
8	44	51.8	433	1	APB_CHICK
9	43	50.6	420	1	DCDA_AQUAE
10	43	50.6	654	1	PGKT_THEMA
11	43	50.6	1075	1	Y124_METJA
12	43	50.6	1120	1	KEFA_ECOLI
13	43	50.6	1166	1	ADDB_BACSU
14	43	50.6	1214	1	BRF1_HUMAN
15	42	49.4	106	1	YBC8_YEAST
16	42	49.4	155	1	YF45_PSEAE
17	42	49.4	218	1	PURQ_METTH
18	42	49.4	307	1	YJM6_YEAST
19	42	49.4	315	1	RPOA_CLOAB
20	42	49.4	322	1	T121_ARCFU
21	42	49.4	393	1	BM15_SHEEP
22	42	49.4	437	1	MYC_FLVTT
23	42	49.4	438	1	MYC_CALJA
24	42	49.4	439	1	MYC_CANFA
25	42	49.4	439	1	MYC_FELCA
26	42	49.4	439	1	MYC_HUMAN
27	42	49.4	439	1	MYC_HYLLA
28	42	49.4	439	1	MYC_MOUSE
29	42	49.4	439	1	MYC_PANTR
30	42	49.4	439	1	MYC_RAT
31	42	49.4	457	1	SYS_SULSO
32	42	49.4	621	1	TRA_STRLI
33	42	49.4	1088	1	RP17_MOUSE
34	42	49.4	1539	1	Y373_HUMAN
35	41	48.2	150	1	TPC2_PONLE
36	41	48.2	226	1	VATE_BOVIN
37	41	48.2	226	1	VATE_HUMAN
38	41	48.2	226	1	YAI5_SCHPO
39	41	48.2	228	1	VATE_MOUSE
40	41	48.2	263	1	MAZG_HAEIN
41	41	48.2	428	1	SYS_PASMU
42	41	48.2	441	1	VATH_ARATH
43	41	48.2	705	1	PPK_BACHD
44	41	48.2	884	1	SYA_RALSO
45	41	48.2	1058	1	CARB_STRPN

P34782 astasia lon
 O34883 bacillus su
 P11682 gallus gall
 O67262 aquifex aeo
 P36204 thermotoga
 Q57588 methanococc
 P77338 escherichia
 P23477 bacillus su
 P55201 homo sapien
 P38202 saccharomyc
 Q04628 pseudomonas
 O26270 methanobact
 P47016 saccharomyc
 Q97ek6 clostridium
 O35003 archaeoglob
 Q9mze2 ovis aries
 P21438 feline leuk
 P49032 callithrix
 Q28350 canis famil
 P06877 felis silve
 P01106 homo sapien
 P49033 hylobates l
 P01108 mus musculu
 P23583 pan troglod
 P09416 rattus norv
 O33780 sulfolobus
 P22409 streptomyce
 Q99nf8 mus musculu
 O15078 homo sapien
 P06708 pontastacus
 P11019 bos taurus
 P36543 homo sapien
 Q09893 schizosacch
 P50518 mus musculu
 P44723 haemophilus
 P57836 pasteurella
 Q9lx65 arabidopsis
 Q9kd27 bacillus ha
 Q8y193 ralstonia s
 Q97qe4 streptococc

ALIGNMENTS

RESULT 1

2NPD_NEUCR

ID 2NPD_NEUCR STANDARD; PRT; 378 AA.

AC Q01284;

DT 01-NOV-1997 (Rel. 35, Created)

DT 01-NOV-1997 (Rel. 35, Last sequence update)

DT 30-MAY-2000 (Rel. 39, Last annotation update)

DE 2-nitropropane dioxygenase precursor (EC 1.13.11.32) (Nitroalkane

DE oxidase) (2-NPD).

GN NCD-2.

OS *Neurospora crassa*.

OC Eukaryota, Fungi; Ascomycota; Pezizomycotina; Sordariomycetes;

OC Sordariales; Sordariaceae; *Neurospora*.

OX NCBI_TaxID=5141;

RN [1]

RP SEQUENCE FROM N. A., AND CHARACTERIZATION.

RC STRAIN=IFO 6067;

RX MEDLINE=98162064; PubMed=9501443;

RA Gorlatova N., Tchorzewski M., Kurihara T., Soda K., Esaki N.;

RT "Purification, characterization, and mechanism of a flavin

RT mononucleotide-dependent 2-nitropropane dioxygenase from *Neurospora*

RT *crassa*.";

RL Appl. Environ. Microbiol. 64:1029-1033(1998).

CC -!- FUNCTION: CATALYZES THE OXIDATION OF NITROALKANES TO PRODUCE
THE

CC CORRESPONDING CARBONYL COMPOUNDS. IT ACTS ON 2-NITROPROPANE
BETTER

CC THAN ON NITROETHANE AND 1-NITROPROPANE, AND ANIONIC FORMS OF

CC NITROALKANES ARE MUCH BETTER SUBSTRATES THAN ARE NEUTRAL
FORMS.

CC -!- CATALYTIC ACTIVITY: 2 2-nitropropane + O(2) = 2 acetone + 2

CC nitrite.

CC -!- COFACTOR: FMN.

CC -!- SUBUNIT: HOMODIMER.

CC -!- SIMILARITY: SOME, TO FAD-DEPENDENT 2-NITROPROPANE DIOXYGENASE.

CC -----

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CC -----

DR EMBL; U22530; AAA64218.1; -.

DR InterPro; IPR004136; 2nprop_dioxygen.
DR InterPro; IPR003009; FMN_enzyme.
DR Pfam; PF03060; NPD; 1.
KW Oxidoreductase; Dioxygenase; Flavoprotein; FMN.
FT PROPEP 1 15 POTENTIAL.
FT CHAIN 16 378 2-NITROPROPANE DIOXYGENASE.
SQ SEQUENCE 378 AA; 39916 MW; E453EB43FD23E441 CRC64;

Query Match 57.6%; Score 49; DB 1; Length 378;
Best Local Similarity 50.0%; Pred. No. 14;
Matches 9; Conservative 5; Mismatches 4; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18
| |::||| || |::|
Db 354 PAGDIVRELREEAKERIK 371
Search completed: February 10, 2003, 09:50:31
Job time : 14 secs

OM protein - protein search, using sw model

Run on: February 10, 2003, 09:49:05 ; Search time 15 Seconds
(without alignments)
115.361 Million cell updates/sec

Title: US-09-865-989-191
Perfect score: 85
Sequence: 1 PVLDLLRELLEELKQKLK 18

Scoring table: BLOSUM62
Gapop 10.0 , Gapext 0.5

Searched 283224 seqs, 96134422 residues

Total number of hits satisfying chosen parameters: 283224

Minimum DB seq length: 0
Maximum DB seq length: 2000000000

Post-processing: Minimum Match 0%
Maximum Match 100%
Listing first 45 summaries

Database : PIR_73:*

1. pir1 *
2. pir2 *
3. pir3 *
4. pir4.*

Pred No is the number of results predicted by chance to have a score greater than or equal to the score of the result being printed, and is derived by analysis of the total score distribution.

SUMMARIES

Result	No.	Score	Match	Length	DB	ID	Description
	1	49	57.6	182	2	E72412	conserved hypothet

2	49	57.6	378	2	T46693	probable 2-nitropr
3	47	55.3	465	2	S03325	transforming prote
4	47	55.3	785	2	S31299	pre-tRNA processin
5	47	55.3	833	2	G83650	DNA gyrase subunit
6	47	55.3	833	2	T46552	DNA topoisomerase
7	46	54.1	322	1	W2WLE	E2 protein - human
8	45	52.9	341	2	C75436	hypothetical prote
9	45	52.9	558	2	S38614	hypothetical prote
10	44	51.8	134	2	D69994	autolytic amidase
11	44	51.8	334	2	D87416	delta-aminolevulin
12	44	51.8	364	2	T29709	probable aspartate
13	44	51.8	433	2	A29626	apolipoprotein B -
14	44	51.8	487	2	D90563	hypothetical prote
15	44	51.8	566	2	E81404	acetolactate synth
16	44	51.8	791	2	H72552	hypothetical prote
17	44	51.8	842	2	AH1432	DNA gyrase chain A
18	44	51.8	842	2	AH1433	DNA gyrase chain A
19	44	51.8	1078	2	T19745	hypothetical prote
20	43.5	51.2	216	2	H71111	hypothetical prote
21	43	50.6	341	2	AG2079	regulatory protein
22	43	50.6	383	2	H75129	probable transamin
23	43	50.6	396	2	T40559	hypothetical coile
24	43	50.6	398	2	F75017	probable glycine C
25	43	50.6	398	2	E71454	probable glycine C
26	43	50.6	405	2	T49534	hypothetical prote
27	43	50.6	420	1	C70404	diaminopimelate de
28	43	50.6	426	2	AD1706	weakly transcripti
29	43	50.6	522	2	T28323	ORF MSV162 probabl
30	43	50.6	571	2	D90157	hypothetical prote
31	43	50.6	612	2	T05331	hypothetical prote
32	43	50.6	654	2	G72344	phosphoglycerate k
33	43	50.6	1120	2	F90693	mechanosensitive c
34	43	50.6	1120	2	B85544	mechanosensitive c
35	43	50.6	1120	2	H64776	probable membrane
36	43	50.6	1163	2	D64315	type I restriction
37	43	50.6	1166	2	A39432	ATP-dependent deox
38	43	50.6	1214	2	JC2069	zinc-finger protei
39	43	50.6	1777	2	AC2088	serine/threonine k
40	42	49.4	97	2	D89792	conserved hypothet
41	42	49.4	106	2	S45762	hypothetical prote
42	42	49.4	155	2	H83453	hypothetical prote
43	42	49.4	156	2	AH0091	probable flagellar
44	42	49.4	186	2	I46085	proto-oncogene - c
45	42	49.4	218	2	E69091	phosphoribosylform

ALIGNMENTS

RESULT 1

E72412

conserved hypothetical protein - *Thermotoga maritima* (strain MSB8)

C;Species: *Thermotoga maritima*

C;Date: 11-Jun-1999 #sequence_revision 11-Jun-1999 #text_change 21-Jul-2000

C;Accession: E72412

R;Nelson, K.E.; Clayton, R.A.; Gill, S.R.; Gwinn, M.L.; Dodson, R.J.; Haft, D.H.; Hickey, E.K.; Peterson, J.D.; Nelson, W.C.; Ketchum, K.A.; McDonald, L.; Utterback, T.R.; Malek, J.A.; Linher, K.D.; Garrett, M.M.; Stewart, A.M.; Cotton, M.D.; Pratt, M.S.; Phillips, C.A.; Richardson, D.; Heidelberg, J.; Sutton, G.G.; Fleischmann, R.D.; White, O.; Salzberg, S.L.; Smith, H.O.; Venter, J.C.; Fraser, C.M.

Nature 399, 323-329, 1999

A;Title Evidence for lateral gene transfer between Archaea and Bacteria from genome sequence of *Thermotoga maritima*.

A;Reference number: A72200; MUID:99287316; PMID:10360571

A;Accession: E72412

A;Status: preliminary

A;Molecule type: DNA

A;Residues 1-182 <ARN>

A;Cross-references: GB:AE001700; GB:AE000512; NID:g4980636; PIDN:AAD35244.1; PID:g4980644; TIGR:TM0151

A;Experimental source: strain MSB8

C;Genetics:

A;Gene: TM0151

Query Match 57.6%; Score 49; DB 2; Length 182;
Best Local Similarity 61.1%; Pred. No. 20;
Matches 11; Conservative 2; Mismatches 5; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18

||:| |:| ||| |||

Db 161 PVVDSRFEILAEKKNK 178

Search completed: February 10, 2003, 09:50:13

Job time : 17 secs

OM protein - protein search, using sw model

Run on: February 10, 2003, 09:49:05 ; Search time 35 Seconds
(without alignments)
68.529 Million cell updates/sec

Title: US-09-865-989-191
Perfect score: 85
Sequence: 1 PVLDLLRELLEELKQKLK 18

Scoring table: BLOSUM62
Gapop 10.0 , Gapext 0.5

Searched: 908470 seqs, 133250620 residues

Total number of hits satisfying chosen parameters: 908470

Minimum DB seq length: 0
Maximum DB seq length: 2000000000

Post-processing: Minimum Match 0%
Maximum Match 100%
Listing first 45 summaries

Database : A_Geneseq_101002:*

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- 2: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1981.DAT *
- 3: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1982.DAT *
- 4: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1983.DAT *
- 5: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1984.DAT *
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- 13: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1992.DAT *
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- 15: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1994.DAT *

16: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1995.DAT.*
 17: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1996.DAT.*
 18: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1997.DAT.*
 19: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1998.DAT.*
 20: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA1999.DAT.*
 21: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA2000.DAT.*
 22: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA2001.DAT.*
 23: /SIDS2/gcgdata/geneseq/geneseqp-embl/AA2002.DAT.*

Pred. No. is the number of results predicted by chance to have a score greater than or equal to the score of the result being printed, and is derived by analysis of the total score distribution.

SUMMARIES

		%					
Result	Query						
No.	Score	Match	Length	DB	ID	Description	

1	85	100.0	18 20	AAY19378		Lecithin:cholester	
2	85	100.0	18 20	AAY19124		Lecithin:cholester	
3	85	100.0	18 20	AAY18870		Lecithin:cholester	
4	85	100.0	18 20	AAY18607		Lecithin:cholester	
5	81	95.3	18 20	AAY19397		Lecithin:cholester	
6	81	95.3	18 20	AAY19401		Lecithin:cholester	
7	81	95.3	18 20	AAY19380		Lecithin:cholester	
8	81	95.3	18 20	AAY19143		Lecithin:cholester	
9	81	95.3	18 20	AAY19147		Lecithin:cholester	
10	81	95.3	18 20	AAY19126		Lecithin:cholester	
11	81	95.3	18 20	AAY18893		Lecithin:cholester	
12	81	95.3	18 20	AAY18889		Lecithin:cholester	
13	81	95.3	18 20	AAY18872		Lecithin:cholester	
14	81	95.3	18 20	AAY18626		Lecithin:cholester	
15	81	95.3	18 20	AAY18630		Lecithin:cholester	
16	81	95.3	18 20	AAY18609		Lecithin:cholester	
17	78	91.8	18 20	AAY19394		Lecithin:cholester	
18	78	91.8	18 20	AAY19379		Lecithin:cholester	
19	78	91.8	18 20	AAY19384		Lecithin:cholester	
20	78	91.8	18 20	AAY19130		Lecithin:cholester	
21	78	91.8	18 20	AAY19140		Lecithin:cholester	
22	78	91.8	18 20	AAY19125		Lecithin:cholester	
23	78	91.8	18 20	AAY18876		Lecithin:cholester	
24	78	91.8	18 20	AAY18886		Lecithin:cholester	
25	78	91.8	18 20	AAY18871		Lecithin:cholester	
26	78	91.8	18 20	AAY18608		Lecithin:cholester	
27	78	91.8	18 20	AAY18613		Lecithin:cholester	

28	78	91.8	18	20	AAAY18623	Lecithin:cholester
29	77	90.6	18	20	AAAY19381	Lecithin:cholester
30	77	90.6	18	20	AAAY19127	Lecithin:cholester
31	77	90.6	18	20	AAAY19137	Lecithin:cholester
32	77	90.6	18	20	AAAY18883	Lecithin:cholester
33	77	90.6	18	20	AAAY18873	Lecithin:cholester
34	77	90.6	18	20	AAAY18610	Lecithin:cholester
35	77	90.6	18	20	AAAY18620	Lecithin:cholester
36	77	90.6	18	20	AAAY19391	Lecithin:cholester
37	76	89.4	18	20	AAAY19383	Lecithin:cholester
38	76	89.4	18	20	AAAY19129	Lecithin:cholester
39	76	89.4	18	20	AAAY18875	Lecithin:cholester
40	76	89.4	18	20	AAAY18612	Lecithin:cholester
41	75	88.2	18	20	AAAY19392	Lecithin:cholester
42	75	88.2	18	20	AAAY19393	Lecithin:cholester
43	75	88.2	18	20	AAAY19402	Lecithin:cholester
44	75	88.2	18	20	AAAY19386	Lecithin:cholester
45	75	88.2	18	20	AAAY19388	Lecithin:cholester

ALIGNMENTS

RESULT 1

AAAY19378

ID AAY19378 standard; Peptide; 18 AA.

XX

AC AAY19378;

XX

DT 14-JUL-1999 (first entry)

XX

DE Lecithin:cholesterol acyltransferase activation exhibiting peptide #191.

XX

KW Apolipoprotein A-I; agonist; dyslipidemic disorder; dyslipidemia;

KW human; lecithin:cholesterol acyltransferase; LCAT; hypercholesterolaemia;

KW cardiovascular disease; atherosclerosis; restenosis; HDL; apoA-I;

KW high density lipoprotein; hypertriglyceridemia; metabolic syndrome,

KW septic shock.

XX

OS Synthetic.

OS Homo sapiens.

XX

PN WO9916459-A1.

XX

PD 08-APR-1999.

XX

PF 28-SEP-1998; 98WO-US20327.

XX

PR 29-SEP-1997; 97US-0940095.

XX

PA (BUTT/) BUTTNER K.

PA (CORN/) CORNUT I.

PA (DASS/) DASSEUX J.

PA (DUFO/) DUFOURCQ J

PA (METZ/) METZ G.

PA (SEKU/) SEKUL R.

XX

PI Buttner K, Cornut I, Dasseux J, Dufourcq J, Metz G;

PI Sekul R;

XX

DR WPI; 1999-277035/23.

XX

PT Peptide agonists of apolipoprotein A-I

XX

PS Example; Page 126; 280pp; English.

XX

CC The present invention describes an agonist (A) of apolipoprotein A-I
CC (apoA-I) which is a 15-29 residue peptide, or analog, that forms an
CC amphipathic alpha-helix in presence of lipids. (A), and their lipid
CC complexes, are used to treat or prevent diseases associated with
CC dyslipidemia, specifically hypercholesterolaemia, cardiovascular
CC disease, atherosclerosis, restenosis, HDL (high density lipoprotein) or
CC apoA-I deficiency; hypertriglyceridemia and metabolic syndrome, also for
CC treating septic shock. When labeled, (A) can also be used diagnostically
CC to measure serum levels of HDL, in particular the HDL subpopulation that
CC is involved in retrograde cholesterol transport, also to image HDL at
CC e.g. atherosclerotic streaks, and to raise antibodies. AAY19188 to
CC AAY19441 represent lecithin:cholesterol acyltransferase (LCAT) activity
CC exhibiting core peptides, which are apoA-I agonists.

XX

SQ Sequence 18 AA;

Query Match 100.0%; Score 85; DB 20; Length 18;
Best Local Similarity 100.0%; Pred. No. 7e-05;
Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKCLK 18

|||||||

Db 1 PVLDLLRELLEELKQKCLK 18

RESULT 2

AAY19124

ID AAY19124 standard; Peptide; 18 AA.

XX

AC AAY19124;

XX

DT 09-JUL-1999 (first entry)

XX

DE Lecithin:cholesterol acyltransferase activation exhibiting peptide #191.

XX

KW Apolipoprotein A-I; agonist; dyslipidemic disorder; dyslipidemia;

KW human; lecithin:cholesterol acyltransferase; LCAT; hypercholesterolaemia;

KW cardiovascular disease; atherosclerosis; restenosis; HDL; apoA-I;

KW high density lipoprotein; hypertriglyceridemia; metabolic syndrome;

KW septic shock.

XX

OS Synthetic.

OS Homo sapiens.

XX

PN WO9916458-A1.

XX

PD 08-APR-1999.

XX

PF 28-SEP-1998; 98WO-US20326.

XX

PR 29-SEP-1997; 97US-0940096.

XX

PA (BUTT/) BUTTNER K.

PA (CORN/) CORNUT I.

PA (DASS/) DASSEUX J.

PA (METZ/) METZ G.

PA (SEKU/) SEKUL R.

XX

PI Buttner K, Cornut I, Dasseux J, Metz G, Sekul R;

XX

DR WPI; 1999-277034/23.

XX

PT Peptide agonists of apolipoprotein A-I

XX

PS Example; Page 117; 254pp; English.

XX

CC The present invention describes an agonist (A) of apolipoprotein A-I

CC (apoA-I) which is a 15-29 residue peptide, or analog, that forms an

CC amphipathic alpha-helix in presence of lipids. (A), and their lipid

CC complexes, are used to treat or prevent diseases associated with

CC dyslipidemia, specifically hypercholesterolaemia, cardiovascular
CC disease, atherosclerosis, restenosis, HDL (high density lipoprotein) or
CC apoA-I deficiency; hypertriglyceridemia and metabolic syndrome, also for
CC treating septic shock. When labeled, (A) can also be used diagnostically
CC to measure serum levels of HDL, in particular the HDL subpopulation that
CC is involved in retrograde cholesterol transport, also to image HDL at
CC e.g. atherosclerotic streaks, and to raise antibodies. AAY18934 to
CC AAY19187 represent lecithin:cholesterol acyltransferase (LCAT) activity
CC exhibiting core peptides, which are apoA-I agonists.

XX

SQ Sequence 18 AA;

Query Match 100.0%; Score 85; DB 20; Length 18;
Best Local Similarity 100.0%; Pred. No. 7e-05;
Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18

|||||||

Db 1 PVLDLLRELLEELKQKLK 18

RESULT 3

AAY18870

ID AAY18870 standard; Peptide; 18 AA.

XX

AC AAY18870;

XX

DT 09-JUL-1999 (first entry)

XX

DE Lecithin:cholesterol acyltransferase activation exhibiting peptide #191.

XX

KW Apolipoprotein A-I; agonist; dyslipidemic disorder; dyslipidemia;

KW human; lecithin:cholesterol acyltransferase; LCAT; hypercholesterolaemia;

KW cardiovascular disease; atherosclerosis; restenosis; HDL; apoA-I;

KW high density lipoprotein; hypertriglyceridemia; metabolic syndrome;

KW septic shock.

XX

OS Synthetic.

OS Homo sapiens.

XX

PN WO9916408-A2.

XX

PD 08-APR-1999.

XX

PF 28-SEP-1998; 98WO-US20328.

XX

PR 29-SEP-1997; 97US-0940093.

XX

PA (BUTT/) BUTTNER K.

PA (CORN/) CORNUT I.

PA (DASS/) DASSEUX J.

PA (METZ/) METZ G.

PA (SEKU/) SEKUL R.

XX

PI Buttner K, Cornut I, Dasseux J, Metz G, Sekul R;

XX

DR WPI; 1999-277031/23.

XX

PT Peptide agonists of apolipoprotein A-I

XX

PS Claim 15; Page 114; 152pp; English.

XX

CC The present invention describes an agonist (A) of apolipoprotein A-I
CC (apoA-I) which is a 14-22 residue peptide, or analog, that forms an
CC amphipathic alpha-helix in presence of lipids. (A), and their lipid
CC complexes, are used to treat or prevent diseases associated with
CC dyslipidemia, specifically hypercholesterolaemia, cardiovascular
CC disease, atherosclerosis, restenosis, HDL (high density lipoprotein) or
CC apoA-I deficiency; hypertriglyceridemia and metabolic syndrome, also for
CC treating septic shock. When labeled, (A) can also be used diagnostically
CC to measure serum levels of HDL, in particular the HDL subpopulation that
CC is involved in retrograde cholesterol transport, also to image HDL at
CC e.g. atherosclerotic streaks, and to raise antibodies. AAY18680 to
CC AAY18933 represent lecithin:cholesterol acyltransferase (LCAT) activity
CC exhibiting core peptides, which are apoA-I agonists.

XX

SQ Sequence 18 AA;

Query Match 100.0%; Score 85; DB 20; Length 18;

Best Local Similarity 100.0%; Pred. No. 7e-05;

Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18

|||||||

Db 1 PVLDLLRELLEELKQKLK 18

RESULT 4

AAY18607

ID AAY18607 standard; Peptide; 18 AA.

XX
AC AAY18607,
XX
DT 09-JUL-1999 (first entry)
XX
DE Lecithin cholesterol acyltransferase activation exhibiting peptide #191.
XX
KW Gene therapy; apolipoprotein A-I; agonist; dyslipidemic disorder;
KW ApoA-I; cardiovascular disease; atherosclerosis; restenosis; LCAT;
KW hyperlipidemia; septic shock; lecithin:cholesterol acyltransferase.
XX
OS Synthetic.
OS Homo sapiens.
XX
PN WO9916409-A2.
XX
PD 08-APR-1999.
XX
PF 28-SEP-1998; 98WO-US20329.
XX
PR 29-SEP-1997; 97US-0940136.
XX
PA (BUTT/) BUTTNER K.
PA (CORN/) CORNUT I.
PA (DASS/) DASSEUX J.
PA (DUFO/) DUFOURCQ J.
PA (METZ/) METZ G.
PA (SEKU/) SEKUL R.
XX
PI Buttner K, Cornut I, Dasseux J, Dufourcq J, Metz G;
PI Sekul R;
XX
DR WPI; 1999-254921/21.
XX
PT Nucleic acid encoding apolipoprotein A-I agonist peptides
XX
PS Claim 49; Page 182, 232pp, English.
XX
CC The present invention describes a nucleic acid (A) encoding an
CC apolipoprotein A-I (apoA-I) agonist (B) that is a peptide, or analog,
CC which forms an amphipathic alpha-helix in presence of lipids. (A),
CC optionally as a complex with lipids, and host cells that contain (A),
CC are useful for gene therapy, or prevention, of diseases associated with
CC dyslipidemia, specifically hypercholesterolaemia, cardiovascular disease,
CC atherosclerosis, restenosis, HDL (high density lipoprotein) and apoA-I

CC deficiency, hypertriglyceridemia and metabolic syndrome, also to treat
CC endotoxemia (septic shock). Host cells containing (A) can also be used
CC to study the role of apoA-I in lipid metabolism. (B) can be used
CC diagnostically, e.g. to measure serum HDL (particularly its
CC subpopulation involved in retrograde cholesterol transport) and for
CC imaging the circulatory system or HDL accumulations at fatty streaks.
CC The present sequence represents a peptide from the present invention.

XX

SQ Sequence 18 AA;

Query Match 100.0%; Score 85; DB 20; Length 18;
Best Local Similarity 100.0%; Pred. No. 7e-05;
Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18

|||||

Db 1 PVLDLLRELLEELKQKLK 18

Search completed: February 10, 2003, 09:49:49

Job time : 36 secs

OM protein - protein search, using sw model

Run on: February 10, 2003, 09:50:35 ; Search time 11 Seconds
(without alignments)
36.277 Million cell updates/sec

Title: US-09-865-989-191
Perfect score 85
Sequence: 1 PVLDLLRELLEELKQKLK 18

Scoring table: BLOSUM62
Gapop 10.0 , Gapext 0.5

Searched: 129505 seqs, 22169297 residues

Total number of hits satisfying chosen parameters: 129505

Minimum DB seq length: 0
Maximum DB seq length: 2000000000

Post-processing: Minimum Match 0%
Maximum Match 100%
Listing first 45 summaries

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3: /cgn2_6/ptodata/2/pubpaa/US06_NEW_PUB.pep.*
4: /cgn2_6/ptodata/2/pubpaa/US06_PUBCOMB.pep.*
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8: /cgn2_6/ptodata/2/pubpaa/US08_PUBCOMB.pep.*
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14: /cgn2_6/ptodata/2/pubpaa/US60_PUBCOMB.pep.*

Pred. No. is the number of results predicted by chance to have a score greater than or equal to the score of the result being printed, and is derived by analysis of the total score distribution.

SUMMARIES

Result No.	% Query					Description
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1	85	100.0	18	9	US-09-865-989-191	Sequence 191, App
2	81	95.3	18	9	US-09-865-989-193	Sequence 193, App
3	81	95.3	18	9	US-09-865-989-210	Sequence 210, App
4	81	95.3	18	9	US-09-865-989-214	Sequence 214, App
5	78	91.8	18	9	US-09-865-989-192	Sequence 192, App
6	78	91.8	18	9	US-09-865-989-197	Sequence 197, App
7	78	91.8	18	9	US-09-865-989-207	Sequence 207, App
8	77	90.6	18	9	US-09-865-989-194	Sequence 194, App
9	77	90.6	18	9	US-09-865-989-204	Sequence 204, App
10	76	89.4	18	9	US-09-865-989-196	Sequence 196, App
11	75	88.2	18	9	US-09-865-989-199	Sequence 199, App
12	75	88.2	18	9	US-09-865-989-201	Sequence 201, App
13	75	88.2	18	9	US-09-865-989-205	Sequence 205, App
14	75	88.2	18	9	US-09-865-989-206	Sequence 206, App
15	75	88.2	18	9	US-09-865-989-215	Sequence 215, App
16	74	87.1	18	9	US-09-865-989-195	Sequence 195, App
17	74	87.1	18	9	US-09-865-989-198	Sequence 198, App
18	74	87.1	18	9	US-09-865-989-200	Sequence 200, App
19	74	87.1	18	9	US-09-865-989-203	Sequence 203, App
20	74	87.1	18	9	US-09-865-989-216	Sequence 216, App
21	74	87.1	18	9	US-09-865-989-221	Sequence 221, App
22	73	85.9	18	9	US-09-865-989-217	Sequence 217, App
23	73	85.9	18	9	US-09-865-989-220	Sequence 220, App
24	73	85.9	18	9	US-09-865-989-227	Sequence 227, App
25	73	85.9	18	9	US-09-865-989-228	Sequence 228, App
26	73	85.9	18	9	US-09-865-989-229	Sequence 229, App
27	73	85.9	18	9	US-09-865-989-231	Sequence 231, App
28	72	84.7	16	9	US-09-865-989-253	Sequence 253, App
29	71	83.5	18	9	US-09-865-989-232	Sequence 232, App
30	69	81.2	18	9	US-09-865-989-209	Sequence 209, App
31	69	81.2	18	9	US-09-865-989-218	Sequence 218, App
32	68	80.0	16	9	US-09-865-989-254	Sequence 254, App
33	68	80.0	18	9	US-09-865-989-219	Sequence 219, App
34	68	80.0	22	9	US-09-865-989-16	Sequence 16, Appl
35	67	78.8	16	9	US-09-865-989-258	Sequence 258, App
36	67	78.8	18	9	US-09-865-989-224	Sequence 224, App

37	67	78.8	22	9	US-09-865-989-130	Sequence 130, App
38	66	77.6	16	9	US-09-865-989-256	Sequence 256, App
39	65	76.5	22	9	US-09-865-989-39	Sequence 39, Appl
40	64	75.3	18	9	US-09-865-989-211	Sequence 211, App
41	64	75.3	22	9	US-09-865-989-1	Sequence 1, Appli
42	64	75.3	22	9	US-09-865-989-3	Sequence 3, Appli
43	64	75.3	22	9	US-09-865-989-4	Sequence 4, Appli
44	64	75.3	22	9	US-09-865-989-6	Sequence 6, Appli
45	64	75.3	22	9	US-09-865-989-8	Sequence 8, Appli

ALIGNMENTS

RESULT 1

US-09-865-989-191

; Sequence 191, Application US/09865989

; Publication No. US20030008827A1

; GENERAL INFORMATION

; APPLICANT: Dasseux, Jean-Louis

; Sekul, Renate

; Buttner, Klaus

; Cornut, Isabelle

; Metz, Gunther

; TITLE OF INVENTION: APOLIPOPROTEIN A-I AGONISTS
; AND THEIR USE TO TREAT DYSLIPIDEMIC DISORDERS

; NUMBER OF SEQUENCES: 258

; CORRESPONDENCE ADDRESS:

; ADDRESSEE: Pennie & Edmonds LLP

; STREET: 1155 Avenue of the Americas

; CITY: New York

; STATE: NY

; COUNTRY: USA

; ZIP: 10036-2811

; COMPUTER READABLE FORM:

; MEDIUM TYPE: Diskette

; COMPUTER: IBM Compatible

; OPERATING SYSTEM: DOS

; SOFTWARE: FastSEQ Version 2.0

; CURRENT APPLICATION DATA:

; APPLICATION NUMBER: US/09/865,989

; FILING DATE: 25-May-2001

; CLASSIFICATION: <Unknown>

; PRIOR APPLICATION DATA:

; APPLICATION NUMBER: 09/465,719

; FILING DATE: 17-DEC-1999
; ATTORNEY/AGENT INFORMATION:
; NAME: Coruzzi, Laura A
; REGISTRATION NUMBER: 30,742
; REFERENCE/DOCKET NUMBER: 009196-0006-999
; TELECOMMUNICATION INFORMATION:
; TELEPHONE: 650-493-4935
; TELEFAX: 650-493-5556
; TELEX: 66141 PENNIE
; INFORMATION FOR SEQ ID NO: 191:
; SEQUENCE CHARACTERISTICS:
; LENGTH: 18 amino acids
; TYPE: amino acid
; STRANDEDNESS: single
; TOPOLOGY: linear
; MOLECULE TYPE: No. US20030008827A1e
; FEATURE:
; NAME/KEY: Other
; LOCATION: 1...18
; OTHER INFORMATION: N-terminal acetylated and
; C-terminal amidated
; SEQUENCE DESCRIPTION: SEQ ID NO: 191:
US-09-865-989-191

Query Match 100.0%; Score 85; DB 9; Length 18;
Best Local Similarity 100.0%; Pred. No. 2.5e-06;
Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18
|||||||
Db 1 PVLDLLRELLEELKQKLK 18

Search completed: February 10, 2003, 09:54:37
Job time : 11 secs

OM protein - protein search, using sw model

Run on: February 10, 2003, 09:49:05 ; Search time 14 Seconds
(without alignments)
37.829 Million cell updates/sec

Title: US-09-865-989-191
Perfect score: 85
Sequence: 1 PVLDLLRELLEELKQKLK 18

Scoring table: BLOSUM62
Gapop 10.0 , Gapext 0.5

Searched: 262574 seqs, 29422922 residues

Total number of hits satisfying chosen parameters: 262574

Minimum DB seq length: 0
Maximum DB seq length: 2000000000

Post-processing: Minimum Match 0%
Maximum Match 100%
Listing first 45 summaries

Database : Issued_Patents_AA.*
1: /cgn2_6/ptodata/1/iaa/5A_COMB.pep:*
2: /cgn2_6/ptodata/1/iaa/5B_COMB.pep:*
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Pred. No. is the number of results predicted by chance to have a score greater than or equal to the score of the result being printed, and is derived by analysis of the total score distribution.

SUMMARIES

Result	%	Query				
No.	Score	Match	Length	DB	ID	Description

1	85	100.0	18	3	US-08-940-095-191	Sequence 191, App
2	85	100.0	18	3	US-08-940-093-191	Sequence 191, App
3	85	100.0	18	3	US-08-940-096-191	Sequence 191, App
4	85	100.0	18	4	US-09-465-719-191	Sequence 191, App
5	85	100.0	18	4	US-09-453-605-191	Sequence 191, App
6	85	100.0	18	4	US-09-453-838-191	Sequence 191, App
7	81	95.3	18	3	US-08-940-095-193	Sequence 193, App
8	81	95.3	18	3	US-08-940-095-210	Sequence 210, App
9	81	95.3	18	3	US-08-940-095-214	Sequence 214, App
10	81	95.3	18	3	US-08-940-093-193	Sequence 193, App
11	81	95.3	18	3	US-08-940-093-210	Sequence 210, App
12	81	95.3	18	3	US-08-940-093-214	Sequence 214, App
13	81	95.3	18	3	US-08-940-096-193	Sequence 193, App
14	81	95.3	18	3	US-08-940-096-210	Sequence 210, App
15	81	95.3	18	3	US-08-940-096-214	Sequence 214, App
16	81	95.3	18	4	US-09-465-719-193	Sequence 193, App
17	81	95.3	18	4	US-09-465-719-210	Sequence 210, App
18	81	95.3	18	4	US-09-465-719-214	Sequence 214, App
19	81	95.3	18	4	US-09-453-605-193	Sequence 193, App
20	81	95.3	18	4	US-09-453-605-210	Sequence 210, App
21	81	95.3	18	4	US-09-453-605-214	Sequence 214, App
22	81	95.3	18	4	US-09-453-838-193	Sequence 193, App
23	81	95.3	18	4	US-09-453-838-210	Sequence 210, App
24	81	95.3	18	4	US-09-453-838-214	Sequence 214, App
25	78	91.8	18	3	US-08-940-095-192	Sequence 192, App
26	78	91.8	18	3	US-08-940-095-197	Sequence 197, App
27	78	91.8	18	3	US-08-940-095-207	Sequence 207, App
28	78	91.8	18	3	US-08-940-093-192	Sequence 192, App
29	78	91.8	18	3	US-08-940-093-197	Sequence 197, App
30	78	91.8	18	3	US-08-940-093-207	Sequence 207, App
31	78	91.8	18	3	US-08-940-096-192	Sequence 192, App
32	78	91.8	18	3	US-08-940-096-197	Sequence 197, App
33	78	91.8	18	3	US-08-940-096-207	Sequence 207, App
34	78	91.8	18	4	US-09-465-719-192	Sequence 192, App
35	78	91.8	18	4	US-09-465-719-197	Sequence 197, App
36	78	91.8	18	4	US-09-465-719-207	Sequence 207, App
37	78	91.8	18	4	US-09-453-605-192	Sequence 192, App
38	78	91.8	18	4	US-09-453-605-197	Sequence 197, App
39	78	91.8	18	4	US-09-453-605-207	Sequence 207, App
40	78	91.8	18	4	US-09-453-838-192	Sequence 192, App
41	78	91.8	18	4	US-09-453-838-197	Sequence 197, App
42	78	91.8	18	4	US-09-453-838-207	Sequence 207, App
43	77	90.6	18	3	US-08-940-095-194	Sequence 194, App
44	77	90.6	18	3	US-08-940-095-204	Sequence 204, App

ALIGNMENTS

RESULT 1

US-08-940-095-191

; Sequence 191, Application US/08940095

; Patent No. 6004925

; GENERAL INFORMATION:

; APPLICANT: Dasseux, Jean-Louis

; APPLICANT: Sekul, Renate

; APPLICANT: Buttner, Klaus

; APPLICANT: Cornut, Isabelle

; APPLICANT: Metz, Gunther

; APPLICANT: Dufourcq, Jean

; TITLE OF INVENTION: APOLIPOPROTEIN A-I AGONISTS

; TITLE OF INVENTION: AND THEIR USE TO TREAT DYSLIPIDEMIC DISORDERS

; NUMBER OF SEQUENCES: 258

; CORRESPONDENCE ADDRESS:

; ADDRESSEE: Pennie & Edmonds LLP

; STREET: 1155 Avenue of the Americas

; CITY: New York

; STATE: NY

; COUNTRY: USA

; ZIP: 10036-2811

; COMPUTER READABLE FORM:

; MEDIUM TYPE: Diskette

; COMPUTER: IBM Compatible

; OPERATING SYSTEM: DOS

; SOFTWARE: FastSEQ Version 2.0

; CURRENT APPLICATION DATA:

; APPLICATION NUMBER: US/08/940,095

; FILING DATE: 29-SEP-1997

; CLASSIFICATION: 514

; PRIOR APPLICATION DATA:

; APPLICATION NUMBER:

; FILING DATE:

; ATTORNEY/AGENT INFORMATION:

; NAME: Coruzzi, Laura A

; REGISTRATION NUMBER: 30,742

; REFERENCE/DOCKET NUMBER: 009196-0004-999

; TELECOMMUNICATION INFORMATION:

; TELEPHONE: 650-493-4935

, TELEFAX: 650-493-5556
, TELEX: 66141 PENNIE
, INFORMATION FOR SEQ ID NO: 191:
, SEQUENCE CHARACTERISTICS:
, LENGTH: 18 amino acids
, TYPE: amino acid
, STRANDEDNESS: single
, TOPOLOGY: linear
, MOLECULE TYPE: No. 6004925e
, FEATURE:
, NAME/KEY: Other
, LOCATION: 1...18
, OTHER INFORMATION: N-terminal acetylated and
, OTHER INFORMATION: C-terminal amidated
US-08-940-095-191

Query Match 100.0%; Score 85; DB 3; Length 18;
Best Local Similarity 100.0%; Pred. No. 1.8e-05;
Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18
|||||||
Db 1 PVLDLLRELLEELKQKLK 18

RESULT 2

US-08-940-093-191
, Sequence 191, Application US/08940093
, Patent No. 6037323
, GENERAL INFORMATION:
, APPLICANT: Dasseux, Jean-Louis
, APPLICANT: Sekul, Renate
, APPLICANT: Buttner, Klaus
, APPLICANT: Cornut, Isabelle
, APPLICANT: Metz, Gunther
, TITLE OF INVENTION: APOLIPOPROTEIN A-I AGONISTS
, TITLE OF INVENTION: AND THEIR USE TO TREAT DYSLIPIDEMIC DISORDERS
, NUMBER OF SEQUENCES: 258
, CORRESPONDENCE ADDRESS:
, ADDRESSEE: Pennie & Edmonds LLP
, STREET: 1155 Avenue of the Americas
, CITY: New York
, STATE: NY
, COUNTRY: USA
, ZIP: 10036-2811

COMPUTER READABLE FORM:
MEDIUM TYPE: Diskette
COMPUTER: IBM Compatible
OPERATING SYSTEM: DOS
SOFTWARE FastSEQ Version 2.0
CURRENT APPLICATION DATA:
APPLICATION NUMBER: US/08/940,093
FILING DATE: 29-SEP-1997
CLASSIFICATION: 435
PRIOR APPLICATION DATA:
APPLICATION NUMBER:
FILING DATE:
ATTORNEY/AGENT INFORMATION:
NAME: Coruzzi, Laura A
REGISTRATION NUMBER: 30,742
REFERENCE/DOCKET NUMBER: 009196-0006-999
TELECOMMUNICATION INFORMATION:
TELEPHONE: 650-493-4935
TELEFAX: 650-493-5556
TELEX: 66141 PENNIE
INFORMATION FOR SEQ ID NO: 191:
SEQUENCE CHARACTERISTICS:
LENGTH 18 amino acids
TYPE: amino acid
STRANDEDNESS: single
TOPOLOGY: linear
MOLECULE TYPE: No. 6037323e
FEATURE
NAME/KEY: Other
LOCATION: 1...18
OTHER INFORMATION: N-terminal acetylated and
OTHER INFORMATION: C-terminal amidated
US-08-940-093-191

Query Match 100.0%, Score 85; DB 3; Length 18;
Best Local Similarity 100.0%; Pred No 1.8e-05;
Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18
|||||||
Db 1 PVLDLLRELLEELKQKLK 18

RESULT 3
US-08-940-096-191

, Sequence 191, Application US/08940096
, Patent No. 6046166
, GENERAL INFORMATION:
, APPLICANT: Dasseux, Jean-Louis
, APPLICANT: Sekul, Renate
, APPLICANT: Buttner, Klaus
, APPLICANT: Cornut, Isabelle
, APPLICANT: Metz, Gunther
, TITLE OF INVENTION: APOLIPOPROTEIN A-I AGONISTS
, TITLE OF INVENTION: AND THEIR USE TO TREAT DYSLIPIDEMIC DISORDERS
, NUMBER OF SEQUENCES: 258
, CORRESPONDENCE ADDRESS:
, ADDRESSEE: Pennie & Edmonds LLP
, STREET: 1155 Avenue of the Americas
, CITY: New York
, STATE: NY
, COUNTRY: USA
, ZIP: 10036-2811
, COMPUTER READABLE FORM:
, MEDIUM TYPE: Diskette
, COMPUTER: IBM Compatible
, OPERATING SYSTEM: DOS
, SOFTWARE: FastSEQ Version 2.0
, CURRENT APPLICATION DATA:
, APPLICATION NUMBER: US/08/940,096
, FILING DATE: 29-SEP-1997
, CLASSIFICATION: 530
, PRIOR APPLICATION DATA:
, APPLICATION NUMBER:
, FILING DATE:
, ATTORNEY/AGENT INFORMATION:
, NAME: Coruzzi, Laura A
, REGISTRATION NUMBER: 30,742
, REFERENCE/DOCKET NUMBER: 009196-0005-999
, TELECOMMUNICATION INFORMATION:
, TELEPHONE: 650-493-4935
, TELEFAX: 650-493-5556
, TELEX: 66141 PENNIE
, INFORMATION FOR SEQ ID NO: 191:
, SEQUENCE CHARACTERISTICS:
, LENGTH: 18 amino acids
, TYPE: amino acid
, STRANDEDNESS: single
, TOPOLOGY: linear
, MOLECULE TYPE: No. 6046166e

; FEATURE:
; NAME/KEY: Other
; LOCATION: 1...18
; OTHER INFORMATION: N-terminal acetylated and
; OTHER INFORMATION: C-terminal amidated
US-08-940-096-191

Query Match 100.0%, Score 85; DB 3; Length 18;
Best Local Similarity 100.0%; Pred. No. 1.8e-05;
Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18
|||||||
Db 1 PVLDLLRELLEELKQKLK 18

RESULT 4

US-09-465-719-191

; Sequence 191, Application US/09465719
; Patent No 6265377

; GENERAL INFORMATION:

; APPLICANT Dasseux, Jean-Louis
; APPLICANT Sekul, Renate
; APPLICANT Buttner, Klaus
; APPLICANT Cornut, Isabelle
; APPLICANT Metz, Gunther

; TITLE OF INVENTION APOLIPOPROTEIN A-I AGONISTS

; TITLE OF INVENTION AND THEIR USE TO TREAT DYSLIPIDEMIC DISORDERS

; NUMBER OF SEQUENCES 258

; CORRESPONDENCE ADDRESS:

; ADDRESSEE: Pennie & Edmonds LLP
; STREET: 1155 Avenue of the Americas
; CITY: New York
; STATE: NY
; COUNTRY USA
; ZIP: 10036-2811

; COMPUTER READABLE FORM:

; MEDIUM TYPE: Diskette
; COMPUTER: IBM Compatible
; OPERATING SYSTEM: DOS
; SOFTWARE: FastSEQ Version 2.0

; CURRENT APPLICATION DATA:

; APPLICATION NUMBER: US/09/465,719
; FILING DATE:
; CLASSIFICATION:

; PRIOR APPLICATION DATA:
 ; APPLICATION NUMBER: 08/940,093
 ; FILING DATE: 29-SEP-1997
 ; ATTORNEY/AGENT INFORMATION:
 ; NAME: Coruzzi, Laura A
 ; REGISTRATION NUMBER: 30,742
 ; REFERENCE/DOCKET NUMBER: 009196-0006-999
 ; TELECOMMUNICATION INFORMATION:
 ; TELEPHONE: 650-493-4935
 ; TELEFAX: 650-493-5556
 ; TELEX: 66141 PENNIE
 ; INFORMATION FOR SEQ ID NO: 191:
 ; SEQUENCE CHARACTERISTICS:
 ; LENGTH: 18 amino acids
 ; TYPE: amino acid
 ; STRANDEDNESS: single
 ; TOPOLOGY: linear
 ; MOLECULE TYPE: No. 6265377e
 ; FEATURE:
 ; NAME/KEY: Other
 ; LOCATION: 1...18
 ; OTHER INFORMATION N-terminal acetylated and
 ; OTHER INFORMATION C-terminal amidated
 US-09-465-719-191

Query Match 100.0%; Score 85; DB 4; Length 18;
 Best Local Similarity 100.0%; Pred. No. 1.8e-05;
 Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18
 |||||
 Db 1 PVLDLLRELLEELKQKLK 18

RESULT 5

US-09-453-605-191

; Sequence 191, Application US/09453605

; Patent No. 6329341

; GENERAL INFORMATION:

; APPLICANT: Dasseux, Jean-Louis
 ; Sekul, Renate
 ; Buttner, Klaus
 ; Cornut, Isabelle
 ; Metz, Gunther
 ; Dufourcq, Jean

TITLE OF INVENTION: APOLIPOPROTEIN A-I AGONISTS
AND THEIR USE TO TREAT DYSLIPIDEMIC DISORDERS

NUMBER OF SEQUENCES: 258

CORRESPONDENCE ADDRESS:

ADDRESSEE: Pennie & Edmonds LLP
STREET: 1155 Avenue of the Americas
CITY: New York
STATE: NY
COUNTRY: USA
ZIP: 10036-2811

COMPUTER READABLE FORM:

MEDIUM TYPE: Diskette
COMPUTER: IBM Compatible
OPERATING SYSTEM: DOS
SOFTWARE: FastSEQ Version 2.0

CURRENT APPLICATION DATA:

APPLICATION NUMBER US/09/453,605
FILING DATE: 26-No. 6329341-1999
CLASSIFICATION: <Unknown>

PRIOR APPLICATION DATA:

APPLICATION NUMBER 08/940,095
FILING DATE: <Unknown>

ATTORNEY/AGENT INFORMATION:

NAME: Coruzzi, Laura A
REGISTRATION NUMBER: 30,742
REFERENCE/DOCKET NUMBER: 009196-0004-999

TELECOMMUNICATION INFORMATION:

TELEPHONE: 650-493-4935
TELEFAX: 650-493-5556
TELEX: 66141 PENNIE

INFORMATION FOR SEQ ID NO 191:

SEQUENCE CHARACTERISTICS:

LENGTH: 18 amino acids
TYPE: amino acid
STRANDEDNESS: single
TOPOLOGY: linear

MOLECULE TYPE: No. 6329341e

FEATURE:

NAME/KEY: Other
LOCATION: 1...18
OTHER INFORMATION: N-terminal acetylated and
C-terminal amidated

SEQUENCE DESCRIPTION: SEQ ID NO: 191:

US-09-453-605-191

Query Match 100.0%; Score 85; DB 4; Length 18;
Best Local Similarity 100.0%; Pred. No. 1.8e-05;
Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKLK 18
|||||||
Db 1 PVLDLLRELLEELKQKLK 18

RESULT 6

US-09-453-838-191

; Sequence 191, Application US/09453838

; Patent No. 6376464

; GENERAL INFORMATION:

; APPLICANT: Dasseux, Jean-Louis

; APPLICANT: Sekul, Renate

; APPLICANT: Buttner, Klaus

; APPLICANT: Cornut, Isabelle

; APPLICANT: Metz, Gunther

; APPLICANT: Dufourcq, Jean

; TITLE OF INVENTION: APOLIPOPROTEIN A-I AGONISTS

; TITLE OF INVENTION: AND THEIR USE TO TREAT DYSLIPIDEMIC DISORDERS

; NUMBER OF SEQUENCES: 258

; CORRESPONDENCE ADDRESS:

; ADDRESSEE: Pennie & Edmonds LLP

; STREET: 1155 Avenue of the Americas

; CITY: New York

; STATE: NY

; COUNTRY: USA

; ZIP: 10036-2811

; COMPUTER READABLE FORM:

; MEDIUM TYPE: Diskette

; COMPUTER: IBM Compatible

; OPERATING SYSTEM: DOS

; SOFTWARE: FastSEQ Version 2.0

; CURRENT APPLICATION DATA:

; APPLICATION NUMBER: US/09/453,838

; FILING DATE:

; CLASSIFICATION:

; PRIOR APPLICATION DATA:

; APPLICATION NUMBER: 08/940,095

; FILING DATE:

; ATTORNEY/AGENT INFORMATION:

; NAME: Coruzzi, Laura A

; REGISTRATION NUMBER: 30,742

REFERENCE/DOCKET NUMBER: 009196-0004-999
TELECOMMUNICATION INFORMATION:
TELEPHONE: 650-493-4935
TELEFAX: 650-493-5556
TELEX: 66141 PENNIE
INFORMATION FOR SEQ ID NO: 191:
SEQUENCE CHARACTERISTICS:
LENGTH: 18 amino acids
TYPE: amino acid
STRANDEDNESS: single
TOPOLOGY: linear
MOLECULE TYPE: No. 6376464e
FEATURE:
NAME/KEY: Other
LOCATION: 1...18
OTHER INFORMATION: N-terminal acetylated and
OTHER INFORMATION: C-terminal amidated
US-09-453-838-191

Query Match 100.0%; Score 85; DB 4; Length 18;
Best Local Similarity 100.0%; Pred. No. 1.8e-05;
Matches 18; Conservative 0; Mismatches 0; Indels 0; Gaps 0;

Qy 1 PVLDLLRELLEELKQKCLK 18
|||||||
Db 1 PVLDLLRELLEELKQKCLK 18

Search completed: February 10, 2003, 09:51:26
Job time : 14 secs